AMENDMENTS TO THE CLAIMS:

This listing of claims will replace all prior versions, and listings, of claims in the application:

(Currently amended) A compound of the formula
 (I):

and pharmaceutically acceptable salts thereof;
wherein:

A is R'-C(O)-, wherein R' is

each R^1 is independently selected from -C(O)-, $-S(O)_2$ -, -C(O)-C(O)-, -O-C(O)-, -O- $S(O)_2$, $-NR^2$ -, $-NR^2$ - $S(O)_2$ -, $-NR^2$ - $S(O)_2$ -, $-NR^2$ - $S(O)_3$ -, $-NR^3$ -, -

each Ht is independently selected from C_3 - C_7 cycloalkyl; C_5 - C_7 cycloalkenyl; C_6 - C_{14} aryl; or a 5-7 membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, $N(R^2)$, O, S and $S(O)_n$; wherein

said aryl or said heterocycle is optionally fused to Q; and wherein any member of said Ht is optionally substituted with one or more substituents independently selected from oxo, -OR², SR^2 , $-\dot{R}^2$, $-N(R^2)(R^2)$, $-R^2$ -OH, -CN, $-CO_2R^2$, $-C(O)-N(R^2)_2$, $-S(O)_2-N(R^2)_2$, $-N(R^2)-C(O)-R^2$, $-N(R^2)-C(O)O-R^2$, $-C(O)-R^2$, $-S(O)_n-R^2$, $-OCF_3$, $-S(O)_n-Q$, methylenedioxy, $-N(R^2)-S(O)_2(R^2)$, halo, $-CF_3$, $-NO_2$, Q, -OQ, $-OR^7$, $-SR^7$, $-R^7$, $-N(R^2)(R^7)$ or $-N(R^7)_2$; each R² is independently selected from H, or C₁-C₄ alkyl optionally substituted with a 3-7 membered saturated, partially saturated or unsaturated carbocyclic ring system; or a 5-7 membered saturated, partially saturated or unsaturated heterocyclic ring containing one or more heteroatoms selected from O, N, S, $S(O)_n$ or $N(R^{33})$; wherein any of said ring systems or $N(R^{33})$ is optionally substituted with 1 to 4 substituents independently selected from -X'-Y', -O-arylalkyl, -S-arylalkyl, $-N(Y')_2$, -N(H)-arylalkyl, $-N(C_1-C_4)$ alkyl)-arylalkyl, oxo, -O-(C_1-C_4 alkyl), OH, C_1-C_4 alkyl, -SO₂H, $-SO_2-(C_1-C_4 \text{ alkyl})$, $-SO_2-NH_2$, $-SO_2-NH(C_1-C_4 \text{ alkyl})$, $-SO_2-N(C_1-C_4)$ $alkyl)_2$, $-NH_2$, $-NH(C_1-C_4 \ alkyl)$, $-N(C_1-C_4 \ alkyl)_2$, -NH-C(O)H, $-N(C_1-C_4 \text{ alkyl})-C(O)H$, $-NH-C(O)-C_1-C_4 \text{ alkyl}$, $-C_1-C_4 \text{ alkyl}-OH$, -OH, -CN, -C(O)OH, $-C(O)O-C_1-C_4$ alkyl, $-C(O)-NH_2$, $-C(0)-NH(C_1-C_4 \text{ alkyl})$, $-C(0)-N(C_1-C_4 \text{ alkyl})_2$, halo or $-CF_3$; X' is -O-, -S-, -NH-, -NHC(O)-, -NHC(O)O-, -NHSO₂-, or $-N(C_1-C_4)$ alkyl-;

Y' is C_1 - C_{15} alkyl, C_2 - C_{15} alkenyl or alkynyl, wherein one to five carbon atoms in Y are optionally substituted with C_3 - C_7 cycloalkyl or C_5 - C_6 cycloalkenyl, C_6 - C_{14} aryl or a 5-7 membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, NH, O, S and S(O)_n;

each R^3 is independently selected from H, Ht, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_3 - C_6 cycloalkyl or C_5 - C_6 cycloalkenyl; wherein any member of said R^3 , except H, is optionally substituted with one or more substituents selected from $-OR^2$, $-C(O)-N(R^2)_2$, $-S(O)_n-N(R^2)_2$, $-N(R^2)_2$, $-N(R^2)-C(O)O(R^2)$, $-N(R^2)-C(O)N(R^2)_2$, $-N(R^2)-C(O)-R^2$; Ht, -CN, $-SR^2$, $-C(O)OR^2$, $N(R^2)-C(O)-R^2$;

each R^{33} is selected from H, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_3 - C_6 cycloalkyl or C_5 - C_6 cycloalkenyl, C_6 - C_{14} aryl or a 5-7 membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, NH, O, S and $S(0)_n$;

each n is independently 1 or 2;

G, when present, is selected from H, R^7 or C_1 - C_4 alkyl, or, when G is C_1 - C_4 alkyl, G and R^7 are bound to one another either directly or through a C_1 - C_3 linker to form a heterocyclic ring; or

when G is not present (i.e., when x in $(G)_x$ is 0), then the nitrogen to which G is attached is bound directly to

the R^7 group in $-OR^7$ with the concomitant displacement of one - ZM group from R^7 ;

D is selected from C_1 - C_6 alkyl which is substituted with Q, which is optionally substituted with one or more groups selected from C_3 - C_6 cycloalkyl, $-R^3$, -O-Q or Q; C_2 - C_4 alkenyl which is substituted with Q, which is optionally substituted with one or more groups selected from $-OR^2$, -S-Ht, $-R^3$, -O-Q or Q; C_3 - C_6 cycloalkyl, which is optionally substituted with or fused to Q; or C_5 - C_6 cycloalkenyl, which is optionally substituted with or fused to Q;

each Q is independently selected from a 3-7 membered saturated, partially saturated or unsaturated carbocyclic ring system; or a 5-7 membered saturated, partially saturated or unsaturated heterocyclic ring containing one or more heteroatoms selected from O, N, S, S(O)_n or N(R²); wherein Q contains one substituent selected from $-OR^2$, $-OR^8$, -O-arylalkyl, $-SR^8$, -S-arylalkyl, $-N(R^2)R^8$ [[,]] or $-N(R^2)$ -arylalkyl and may be optionally substituted with one or more additional substituents independently selected from oxo, $-OR^8$, -O-arylalkyl $-SR^8$, -S-arylalkyl, $-N(R^2)R^8$, $-N(R^2)$ -arylalkyl, $-OR^2$, $-R^2$, $-SO_2R^2$, $-SO_2-N(R^2)_2$, $-N(R^2)_2$, $-N(R^2)$ -arylalkyl, $-OR^2$, $-R^2$, $-SO_2R^2$, $-SO_2-N(R^2)_2$, $-N(R^2)_2$, halo or $-CF_3$;

each R^8 is independently selected from Ht, $-C_1-C_{15}$

branched or straight chain alkyl, alkenyl or alkynyl wherein one to five carbon atoms in said alkyl, alkenyl or alkynyl are independently replaced by W, or wherein one to five carbon atoms in said alkyl, alkenyl or alkynyl are substituted with Ht; and wherein R^8 is additionally and optionally substituted with one or more groups independently selected from -OH, -S(C₁-C₆ alkyl), -CN, -CF₃, -N(R^2)₂, halo, -C₁-C₄-alkyl, -C₁-C₄-alkoxy; -Ht; -O-Ht; -NR²-CO-N(R^2)₂; -CO-N(R^2)₂; -R¹-C₂-C₆ alkenyl, which is optionally substituted with one or more groups independently selected from hydroxy, C₁-C₄ alkoxy, Ht, -O-Ht, -NR²-CO-N(R^2)₂ or -CO-N(R^2)₂; or R^7 ;

wherein W is -O-, $-NR^2$ -, -S-, -C(O)-, -C(S)-, $-C(=NR^2)$ -, $-S(O)_2$ -, $-NR^2$ - $S(O)_2$ -, $-S(O)_2$ -, $-S(O)_2$ -, $-NR^2$ -C(O)-, -O-C(O)-, -O-C(O)-, -O-C(O)-, -O-C(O)-, -O-C(O)-, -O-, -O-

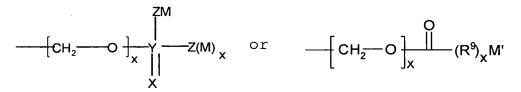
D' is selected from C_1 - C_{15} alkyl, C_1 - C_{15} alkoxy, C_2 - C_{15} alkenyl, C_2 - C_{15} alkenyloxy, C_2 - C_{15} alkynyl, or C_2 - C_{15} alkynyloxy, wherein D' optionally comprises one or more substituents independently selected from Ht, oxo, halo, -CF₃, -OCF₃, -NO₂, azido, -SH, -SR³, -N(R³)-N(R³)₂, -O-N(R³)₂, -(R³)N-O-(R³), -N(R³)₂, -CN, -CO₂R³, -C(O)-N(R³)₂, -S(O)_n-N(R³)₂, -N(R³)-C(O)-R³, -N(R³)-C(O)-N(R³)₂, -C(O)-R³, -S(O)_n-R³, -N(R³)-S(O)_n(R³), -N(R³)-S(O)_n-N(R³)₂, -S-NR³-C(O)R³, -C(S)N(R³)₂, -C(S)R³, -NR³-C(O)OR³, -O-C(O)OR³, -O-C(O)N(R³)₂, -NR³-C(S)R³, =N-OH,

 $= N - OR^{3}, = N - N(R^{3})_{2}, = NR^{3}, = NNR^{3}C(O)N(R^{3})_{2}, = NNR^{3}C(O)OR^{3},$ $= NNR^{3}S(O)_{n} - N(R^{3})_{2}, -NR^{3} - C(S)OR^{3}, -NR^{3} - C(S)N(R^{3})_{2},$ $- NR^{3} - C[=N(R^{3})] - N(R^{3})_{2}, - N(R^{3}) - C[=N - NO_{2}] - N(R^{3})_{2},$ $- N(R^{3}) - C[=N - NO_{2}] - OR^{3}, - OC(O)R^{3}, - OC(S)R^{3}, - OC(O)N(R^{3})_{2},$ $- C(O)N(R^{3}) - N(R^{3})_{2}, - N(R^{3}) - N(R^{3})C(O)R^{3}, - N(R^{3}) - OC(O)R^{3},$ $- N(R^{3}) - OC(O)R^{3}, - N(R^{3}) - OC(O)R^{3}, - OC(S)N(R^{3})_{2}, - OC(S)N(R^{3})(R^{3}),$ $- PO_{3} - R^{3};$

E is selected from Ht; O-Ht; Ht-Ht; Ht fused with Ht; $-O-R^3$; $-N(R^2)(R^3)$; $-N(R^2)-Ht$; C_1-C_6 alkyl, which is optionally substituted with one or more groups selected from R^4 or Ht; C_2-C_6 alkenyl, which is optionally substituted with one or more groups selected from R^4 or Ht; C_3-C_6 saturated carbocycle, which is optionally substituted with one or more groups selected from R^4 or Ht; or C_5-C_6 unsaturated carbocycle, which is optionally substituted with one or more groups selected from R^4 or Ht;

each R^4 is independently selected from $-R^2$, $-OR^2$, $-OR^3$, $-SR^2$, $-SOR^2$, $-SO_2R^2$, $-CO_2R^2$, $-OC(O)-R^2$, $-C(O)-N(R^2)_2$, $-C(O)-NR^2(OR^2)$, $-S(O)_2-N(R^2)_2$, halo, $-NR^2-C(O)-R^2$, $-NR^2-OR^2$, $-N(R^2)_2$ or -CN;

each R⁷ is independently selected from hydrogen,



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wherein each M is independently selected from H, Li, Na, K, Mg, Ca, Ba, $-N(R^2)_4$, C_1-C_{12} -alkyl, C_2-C_{12} -alkenyl, or $-R^6$; wherein 1 to 4 $-CH_2$ radicals of the alkyl or alkenyl group, other than the $-CH_2$ that is bound to Z, is optionally replaced by a heteroatom group selected from O, S, S(O), $S(O_2)$, or $N(R^2)$; and wherein any hydrogen in said alkyl, alkenyl or R^6 is optionally replaced with a substituent selected from $O(C_1-C_4)$ alkyl, $O(C_2-C_4)$ alkyl, $O(C_2-C_4)$ alkyl, $O(C_3-C_4)$ alkyl, $O(C_4-C_4)$ alkyl, $O(C_4-C_4)$

M' is H, C_1 - C_{12} -alkyl, C_2 - C_{12} -alkenyl, or $-R^6$; wherein 1 to 4 - CH_2 radicals of the alkyl or alkenyl group is optionally replaced by a heteroatom group selected from O, S, S(O), $S(O_2)$, or $N(R^2)$; and wherein any hydrogen in said alkyl, alkenyl or R^6 is optionally replaced with a substituent selected from $O(R^2)$, $O(R^2)$

x is 0 or 1;

Z is O, S, $N(R^2)_2$, or, when M is not present, H;

Y is P or S;

X is O or S; and

 R^9 is $C(R^2)_2$, O or $N(R^2)$; and wherein when Y is S, Z is not S; and

 R^6 is a 5-6 membered saturated, partially saturated or unsaturated carbocyclic or heterocyclic ring system, or an 8-10 membered saturated, partially saturated or unsaturated bicyclic ring system; wherein any of said heterocyclic ring systems contains one or more heteroatoms selected from O, N, S, $S(O)_n$ or $N(R^2)$; and wherein any of said ring systems optionally contains 1 to 4 substituents independently selected from -OH, -C₁-C₄ alkyl, -O-(C₁-C₄ alkyl) or -O-C(O)-(C₁-C₄ alkyl).

2. (Original) The compound according to claim 1, wherein R^8 is $-C_1-C_4$ -branched or straight chain alkyl, wherein one to two carbon atoms in said alkyl are independently replaced by W, wherein R^8 is additionally and optionally substituted with one or more groups independently selected from -OH; $-C_1-C_4$ -alkoxy; -Ht; -O-Ht; $-NR^2-CO-N(R^2)_2$; $-CO-N(R^2)_2$; $-R^1-C_2-C_6$ alkenyl, which is optionally substituted with one or more groups independently selected from hydroxy, C_1-C_4 alkoxy, $-C_4$ alkoxy, -

wherein W is -O-, $-NR^2$ -, $-NR^2$ -S(O)₂-, $-NR^2$ -C(O)O-, -O-C(O)NR²-, $-NR^2$ -C(O)NR²-, $-NR^2$ -C(S)NR²-, $-NR^2$ C(O)-, -C(=NR²)-, -C(O)NR²-, $-NR^2$ -C(=N-CN)-NR²-, $-NR^2$ C(=N-CN)O- or -C(O)O-; and

wherein Ht, R^1 , R^2 and R^7 are as defined in claim 1.

3. (Previously presented) The compound according to claim 1, wherein R^8 is a $-C_1-C_4$ -branched or straight alkyl chain, wherein one to two carbon atoms are substituted with Ht;

wherein Ht is C_{6-14} aryl or a 5-7 membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, $N(R^2)$, O, S and $S(O)_n$, wherein any member of Ht is optionally substituted with one or more substituents independently selected from oxo, $-OR^2$, SR^2 , $-R^2$, $-N(R^2)(R^2)$, $-R^2$ -OH, -CN, $-CO_2R^2$, $-C(O)-N(R^2)_2$, $-S(O)_2-N(R^2)_2$, $-N(R^2)-C(O)-R^2$, $-N(R^2)-C(O)-R^2$, $-S(O)_n-R^2$, $-OCF_3$, $-S(O)_n-Q$, methylenedioxy, $-N(R^2)-S(O)_2(R^2)$, halo, $-CF_3$, $-NO_2$, Q, -OQ, $-OR^7$, $-SR^7$, $-R^7$, $-N(R^2)(R^7)$ or $-N(R^7)_2$.

4. (Currently amended) The compound according to claim 1, wherein R^8 is selected from:

5. (Original) The compound according to claim 1, wherein at least one \mathbb{R}^7 is selected from:

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acetyl, $\stackrel{\text{O}}{\smile}$, $\stackrel{\text{O}}{\smile}$, $\stackrel{\text{C}}{\smile}$, $\stackrel{\text{C}}{\smile}$, $\stackrel{\text{C}}{\smile}$, $\stackrel{\text{C}}{\smile}$, $\stackrel{\text{C}}{\smile}$

-(L)-aspartic acid, -(L)-
$$\gamma$$
-t-butyl-aspartic acid,

-(L)-(L)-3-pyridylalanine, -(L)-histidine, -CHO,
$$CF_3$$
,

 PO_3 -spermine, PO_3 -(spermidine)₂ or PO_3 -(meglamine)₂.

6. (Canceled).

7. (Previously presented) The compound according to claim 1, wherein:

D' is $-CH_2-R'$ ', wherein R'' is selected from: isobutyl,;

wherein m is 0 to 3.

8. (Previously presented) The compound according to claim 1, wherein:

E is selected from:

9. (Currently amended) A compound having the

formula (II):

and pharmaceutically acceptable salts thereof; wherein:

A is selected from R'-C(O)-, wherein R' is

each R^1 is independently selected from -C(0)-, $-S(0)_2$ -, -C(0)-C(0)-, -O-C(0)-, $-O-S(0)_2$, $-NR^2$ -, $-NR^2$ - $S(0)_2$ -, $-NR^2$ -S

each Ht is independently selected from C_3 - C_7 cycloalkyl; C_5 - C_7 cycloalkenyl; C_6 - C_{14} aryl; or a 5-7 membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, $N(R^2)$, O, S and $S(O)_n$; wherein said aryl or said heterocycle is optionally fused to Q; and wherein any member of said Ht is optionally substituted with one or more substituents independently selected from oxo, $-OR^2$,

$$\begin{split} & SR^2\,, \ -R^2\,, \ -N\,(R^2)\,(R^2)\,, \ -R^2-OH\,, \ -CN\,, \ -CO_2R^2\,, \ -C\,(O)\,-N\,(R^2)_{\,2}\,, \\ & -S\,(O)_{\,2}-N\,(R^2)_{\,2}\,, \ -N\,(R^2)\,-C\,(O)\,-R^2\,, \ -N\,(R^2)\,-C\,(O)\,O-R^2\,, \ -C\,(O)\,-R^2\,, \\ & -S\,(O)_{\,n}-R^2\,, \ -OCF_3\,, \ -S\,(O)_{\,n}-Q\,, \ \text{methylenedioxy}\,, \ -N\,(R^2)\,-S\,(O)_{\,2}\,(R^2)\,, \\ & \text{halo}\,, \ -CF_3\,, \ -NO_2\,, \ Q\,, \ -OQ\,, \ -OR^7\,, \ -SR^7\,, \ -R^7\,, \ -N\,(R^2)\,(R^7)\,\,\text{or}\,\, -N\,(R^7)_{\,2}\,; \end{split}$$

each R^2 is independently selected from H, or C_1 - C_4 alkyl optionally substituted with a 3-7 membered saturated, partially saturated or unsaturated carbocyclic ring system; or a 5-7 membered saturated, partially saturated or unsaturated heterocyclic ring containing one or more heteroatoms selected from O, N, S, S(O)_n or N(R^{33}); wherein any of said ring systems or N(R^{33}) is optionally substituted with 1 to 4 substituents independently selected from -X'-Y', -O-arylalkyl,

 $-S-arylalkyl, -N(Y')_2, -N(H)-arylalkyl, -N(C_1-C_4) \\ alkyl)-arylalkyl, oxo, -O-(C_1-C_4) alkyl), OH, C_1-C_4 alkyl, -SO_2H, \\ -SO_2-(C_1-C_4) alkyl), -SO_2-NH_2, -SO_2-NH(C_1-C_4) alkyl), -SO_2-N(C_1-C_4) \\ alkyl)_2, -NH_2, -NH(C_1-C_4) alkyl), -N(C_1-C_4) alkyl)_2, -NH-C(O)H, \\ -N(C_1-C_4) alkyl)-C(O)H, -NH-C(O)-C_1-C_4) alkyl, -C_1-C_4 alkyl-OH, \\ -OH, -CN, -C(O)OH, -C(O)O-C_1-C_4) alkyl, -C(O)-NH_2, \\ \\ -OH, -CN, -C(O)OH, -C(O)O-C_1-C_4) alkyl, -C(O)O-C_1-C_4) alkyl,$

 $\label{eq:condition} X' \mbox{ is -O-, -S-, -NH-, -NHC(O)-, -NHC(O)O-, -NHSO_2-,} \\ or -N(C_1-C_4) \mbox{ alkyl-;}$

 $-C(0)-NH(C_1-C_4 \text{ alkyl})$, $-C(0)-N(C_1-C_4 \text{ alkyl})_2$, halo or $-CF_3$;

Y' is C_1 - C_{15} alkyl, C_2 - C_{15} alkenyl or alkynyl, wherein one to five carbon atoms in Y are optionally substituted with C_3 - C_7 cycloalkyl or C_5 - C_6 cycloalkenyl, C_6 - C_{14} aryl or a 5-7

membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, NH, O, S and $S(O)_n$;

each R^3 is independently selected from H, Ht, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_3 - C_6 cycloalkyl or C_5 - C_6 cycloalkenyl; wherein any member of said R^3 , except H, is optionally substituted with one or more substituents selected from $-OR^2$, $-C(O)-N(R^2)_2$, $-S(O)_n-N(R^2)_2$, $-N(R^2)_2$, $-N(R^2)-C(O)O(R^2)_2$, $-N(R^2)-C(O)N(R^2)_2$, $-N(R^2)-C(O)-R^2$; Ht, -CN, $-SR^2$, $-C(O)OR^2$, $N(R^2)-C(O)-R^2$;

each R^{33} is selected from H, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_3 - C_6 cycloalkyl or C_5 - C_6 cycloalkenyl, C_6 - C_{14} aryl or a 5-7 membered saturated or unsaturated heterocycle, containing one or more heteroatoms selected from N, NH, O, S and $S(O)_n$;

each n is independently 1 or 2;

each Q is independently selected from a 3-7 membered saturated, partially saturated or unsaturated carbocyclic ring system; or a 5-7 membered saturated, partially saturated or unsaturated heterocyclic ring containing one or more heteroatoms selected from O, N, S, $S(O)_n$ or $N(R^2)$; wherein Q contains one substituent selected from $-OR^2$, $-OR^8$, -O-arylalkyl, $-SR^8$, -S-arylalkyl, $-N(R^2)R^8$ [[,]] or $-N(R^2)$ -arylalkyl and may be optionally substituted with one or more additional substituents independently selected from oxo,

 $-\mathrm{OR}^8, \ -\mathrm{O-arylalkyl} \ -\mathrm{SR}^8, \ -\mathrm{S-arylalkyl}, \ -\mathrm{N}(R^2)R^8, \ -\mathrm{N}(R^2) -\mathrm{arylalkyl}, \ -\mathrm{OR}^2, \ -\mathrm{R}^2, \ -\mathrm{SO}_2R^2, \ -\mathrm{SO}_2-\mathrm{N}(R^2)_2, \ -\mathrm{N}(R^2)_2, \ -\mathrm{N}(R^2)_2, \ -\mathrm{N}(R^2)_2, \ -\mathrm{OH}, \ (C_1-C_4) -\mathrm{OH}, \ -\mathrm{CN}, \ -\mathrm{CO}_2R^2, \ -\mathrm{C}(\mathrm{O}) -\mathrm{N}(R^2)_2, \ \mathrm{halo}$ or $-\mathrm{CF}_3;$

each R^8 is independently selected from Ht, $-C_1-C_{15}$ branched or straight chain alkyl, alkenyl or alkynyl wherein one to five carbon atoms in said alkyl, alkenyl or alkynyl are independently replaced by W, or wherein one to five carbon atoms in said alkyl, alkenyl or alkynyl are substituted with Ht; and wherein R^8 is additionally and optionally substituted with one or more groups independently selected from -OH, $-S(C_1-C_6$ alkyl), -CN, $-CF_3$, $-N(R^2)_2$, halo, $-C_1-C_4$ -alkyl, $-C_1-C_4$ -alkoxy; -Ht; -O-Ht; $-NR^2-CO-N(R^2)_2$; $-CO-N(R^2)_2$; $-R^1-C_2-C_6$ alkenyl, which is optionally substituted with one or more groups independently selected from hydroxy, C_1-C_4 alkoxy, Ht, -O-Ht, $-NR^2-CO-N(R^2)_2$ or $-CO-N(R^2)_2$; or R^7 ;

wherein W is -O-, -NR²-, -S-, -C(O)-, -C(S)-, -C(=NR²)-, -S(O)₂-, -NR²-S(O)₂-, -S(O)₂-, -NR²-C(O)O-, -O-C(O)NR²-, -NR²-C(O)NR²-, -NR²-C(S)NR²-, -CONR², -NR²C(O)-, -C(S)NR², -NR²C(S)-, -NR²-C(=N-CN)-NR²-, -NR²C(=N-CN)O- or -C(O)O-;

D' is selected from C_1 - C_{15} alkyl, C_1 - C_{15} alkoxy, C_2 - C_{15} alkenyl, C_2 - C_{15} alkenyloxy, C_2 - C_{15} alkynyl, or C_2 - C_{15} alkynyloxy, wherein D' optionally comprises one or more substituents independently selected from Ht, oxo, halo, -CF₃, -OCF₃, -NO₂,

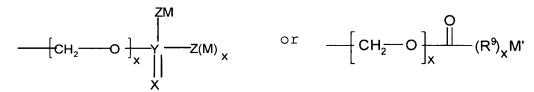
azido, -SH, $-SR^3$, $-N(R^3) - N(R^3)_2$, $-O-N(R^3)_2$, $-(R^3)N-O-(R^3)$, $-N(R^3)_2$, -CN, $-CO_2R^3$, $-C(O) - N(R^3)_2$, $-S(O)_n - N(R^3)_2$, $-N(R^3) - C(O) - R^3$, $-N(R^3) - C(O) - N(R^3)_2$, $-C(O) - R^3$, $-S(O)_n - R^3$, $-N(R^3) - S(O)_n (R^3)$, $-N(R^3) - S(O)_n - N(R^3)_2$, $-S-NR^3 - C(O)R^3$, $-C(S)N(R^3)_2$, $-C(S)R^3$, $-NR^3 - C(O)OR^3$, $-O-C(O)OR^3$, $-O-C(O)N(R^3)_2$, $-NR^3 - C(S)R^3$, -N-OH, $-N-OR^3$, $-N-N(R^3)_2$, $-NR^3 - C(S)OR^3$, -N-OH, $-N-OR^3$, $-N-N(R^3)_2$, $-NR^3 - C(S)OR^3$, $-NR^3 - C(S)N(R^3)_2$, $-NR^3 - C(S)OR^3$, $-NR^3 - C(S)N(R^3)_2$, $-N(R^3) - C(S)OR^3$, $-OC(O)N(R^3)_2$, $-N(R^3) - OC(O)R^3$, $-OC(O)N(R^3)_2$, $-N(R^3)_2$, $-N(R^3)_2$, $-N(R^3)_2$, $-N(R^3)_2$, $-N(R^3)_2$, $-OC(S)N(R^3)_2$, $-OC(S)N(R^3)_$

E is selected from Ht; O-Ht; Ht-Ht; Ht fused with Ht; $-O-R^3$; $-N(R^2)(R^3)$; $-N(R^2)-Ht$; C_1-C_6 alkyl, which is optionally substituted with one or more groups selected from R^4 or Ht; C_2-C_6 alkenyl, which is optionally substituted with one or more groups selected from R^4 or Ht; C_3-C_6 saturated carbocycle, which is optionally substituted with one or more groups selected from R^4 or Ht; or C_5-C_6 unsaturated carbocycle, which is optionally substituted with one or more groups selected from R^4 or Ht;

each R^4 is independently selected from $-R^2$, $-OR^2$, $-OR^3$, $-SR^2$, $-SOR^2$, $-SO_2R^2$, $-CO_2R^2$, $-OC(O) - R^2$, $-C(O) - N(R^2)_2$, $-C(O) - NR^2(OR^2)$, $-S(O)_2 - N(R^2)_2$, halo, $-NR^2 - C(O) - R^2$, $-NR^2 - OR^2$,

 $-N(R^2)_2$ or -CN;

each R⁷ is independently selected from hydrogen,



M' is H, C_1-C_{12} -alkyl, C_2-C_{12} -alkenyl, or $-R^6$; wherein 1 to 4 -CH₂ radicals of the alkyl or alkenyl group is optionally replaced by a heteroatom group selected from O, S, S(O), $S(O_2)$, or $N(R^2)$; and wherein any hydrogen in said alkyl, alkenyl or R^6 is optionally replaced with a substituent selected from oxo, $-OR^2$, $-C_1-C_4$ alkyl, $-N(R^2)_2$, $N(R^2)_3$, -OH, $-O-(C_1-C_4$ alkyl), -CN, $-C(O)OR^2$, $-C(O)-N(R^2)_2$, $-S(O)_2-N(R^2)_2$,

 $-N(R^2)-C(0)-R_2$, $-C(0)R^2$, $-S(0)_n-R^2$, $-OCF_3$, $-S(0)_n-R^6$, $-N(R^2)-S(0)_2(R^2)$, halo, $-CF_3$, or $-NO_2$;

x is 0 or 1;

Z is O, S, $N(R^2)_2$, or, when M is not present, H;

Y is P or S;

X is O or S; and

 $\mbox{\sc R}^9$ is $\mbox{\sc C}(\mbox{\sc R}^2)_2,$ O or $\mbox{\sc N}(\mbox{\sc R}^2)_{\sc z}$ and wherein when Y is S, Z is not S; and

 R^6 is a 5-6 membered saturated, partially saturated or unsaturated carbocyclic or heterocyclic ring system, or an 8-10 membered saturated, partially saturated or unsaturated bicyclic ring system; wherein any of said heterocyclic ring systems contains one or more heteroatoms selected from O, N, S, $S(O)_n$ or $N(R^2)$; and wherein any of said ring systems optionally contains 1 to 4 substituents independently selected from -OH, -C₁-C₄ alkyl, -O-(C₁-C₄ alkyl) or -O-C(O)-(C₁-C₄ alkyl).

10. (Original) The compound according to claim 9, wherein \mathbb{R}^8 is selected from:

11. (Currently amended) The compound according to claim 9, wherein \mathbb{R}^8 is selected from:

12. (Original) The compound according to claim 9,

wherein R⁸ is selected from:

$$O_2N$$
 O_2
 O_2

13. (Original) The compound according to claim 9, wherein $\ensuremath{\text{R}}^8$ is selected from:

14. (Previously presented) The compound according to claim 9, wherein \mathbb{R}^8 is selected from:

15. (Currently amended) The compound according to claim 9, wherein said compound is selected from compound numbers: 26, 27, 31, 33, 35, 36, 38, 41, 43, 48, 49, 51, 52, 53, 54, 55, 56, 57, 59, 60, 71, 72, 73, 74, 202, 203, 209, 213, 215, 223, 227, 231, 233, 236, 237, 239, 243, 247, 250, 260, 263, 271, 281, 289, 293, 295, 309, 317, 319, 320, 322, 334, 335, 348, 364, 367, 368, 375, 382, 383 and 396, wherein said compound numbers are as defined below: [[.]]

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	A	S(O) ₂ E	R ⁸
33	H H	NH ₂	NO
35	H	NH ₂	N
36	H H	SONH2	N
38	H H	SO ₂ OH	
41	H	SO ₂ OH	
43	H H	SO ₂ OH	N
48	H H	- § 2	
49	H H	-S ₂	

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	A	S(O) ₂ E	R ⁸
51	H H	S _{O2}	
52	H H	S _{O2}	
53	H O O O O O O O O O O O O O O O O O O O	S _{O2}	N
54	H 0 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	S ₂	N
55	H	S ₂	NO
56	H H	S _{O2}	N N
57	H H	S _{O2}	N N
59	H	SO ₂ NH ₂	N

A
$$S(0)_{2}E$$
 R^{8}
 71
 R^{9}
 R^{9}

A	S(O) ₂ E	R ⁸
74 H	NH ₂	NO

	A	R ⁸	D,	E
202	H O H		*<	XX°>
203	H O N	*D	×<	↓↓\cdots
209	H O J	s	×. <	
213	H, O, J,	s N	×. <	
215	H, O, J,	X CN	×<	XX°
223	H°,	NHCOO-t-bu	×.<	(C)

		Y	T	γ ··· ··· ··· ··· ··· ··· ··· ··· ··· ·
227	H O N	NHSO₂Me	×.<	
231	H ° ,	× Nº	×. <	
233	H°,	× Y	×.<	(X)
236	HON	÷ Ho	ж. <	\(\sigma_{\circ}\)
237	H°,	→ NH S S S	×. <	
239	H O J		×. <	X°>
243	H ° ,	N O OMe	×. <	XX°
247	H° N	, N O OEt	×. <	III »
250	H°,	H O O-i-Pr	×. <	
260	H°,	NHCSNHMe	×.<	
263	H°,	N.CN N. OPh	×.<	
271	H O J	× H N	×.<	
281	H O J	× NNN	×.<	II.
289	H O O	, D-cn	×.<	NH ₂

r				
293	H	N	×. <	NH ₂
295	H O O	, In	×.<	OH Me
309	H, O, J,	-CONHMe	×. <	(C)
317	H°,	,,,,он ,,,,,он	×. <	\(\sigma\)
319	H° N	ЖОН	×<	(八)
320	H ° ,	OCONHMe	×	XX°
322	H O A	<i>;</i> ОН	×.<	\(\sigma\)
334	H O N	O ₂ N COOH	×. <	↓ǰ>
335	H ° ,	NC (N	×. <	\(\frac{1}{2}\)
348	H ° ,	, L s	×. <	OH
364	H O J	÷0	×. <	\(\frac{\chi}{\chi}\)
367	H O J	; Ocn	×.<	\(\sigma_{\circ}^{\circ}\)
368	H O J	× In	×. <	\(\sigma_{\circ}\)

375	H, O, J,	, NO	×-<	XX.
382	H O J	; CN	×<	
383	H O J	NO ₂	×<	
396	H O J	CN	×.	XX°>

16. (Currently amended) The compound according to claim 15, wherein said compound is selected from compound numbers: 26, 27, 31, 33, 35, 36, 38, 41, 43, 48, 49, 51, 52, 53, 54, 55, 56, 57, 59, 60, 71, 72, 73, 74, 209, 215, 227, 233, 237, 281, 289, 295, 309, 322, 335, 364, 368, 382 and 383, wherein said compound numbers are as defined below: [[.]]

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	A	S(O) ₂ E	R ⁸
26	H H	NH ₂	
27	H H	NH ₂	N O
31	H H	NH ₂	N
33	H H	NH ₂	NO
35	H H	NH ₂	NO
36	H H	S NH ₂	N
38	H H	SO ₂ OH	
41	H H	SO ₂ OH	

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43	H H	SO ₂ OH	N
48	H H	S _{O2}	
49	H H	S _{O₂}	N N
51	H H	-S ₂	N O
52	H H	- S ₂	N
53	H H	S C C C C C C C C C C C C C C C C C C C	N
54	H H	S C C C C C C C C C C C C C C C C C C C	N
55	H	S C C C C C C C C C C C C C C C C C C C	NO
56	H	S C C C C C C C C C C C C C C C C C C C	N N

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A
$$S(0)_2E$$
 R^8

71 H_0 S_0 NH_2 NH_2 S_0 S_0

	A	S(O) ₂ E	R ⁸
73	H H	-S _{O₂}	
74	H	NH ₂	

	A	R ⁸	D'	E
209	H, O, J,	s	×<	
215	H, O, J,	CN	×<	\(\sigma_{\circ}\)
227	H°, ,	NHSO ₂ Me	×	\(\sigma\)
233	H°,	× HO	×	(C)
237	H°,	~ H ~ S ~ S ~ S ~ S ~ S ~ S ~ S ~ S ~ S	×. <	\(\sqrt{\chi}\)

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281	H O J	X N N N	×<	
289	H O O	; D-cn	×	NH ₂
295	H O O	; Sn	×	OH Me
309	H O J	-CONHMe	ж<	\(\sigma\)
322	H O J	; ∕ он	×	
335	H O H	NC N	×	XX°
364	H O N		×<	
368	H, O, I,	× S	\times	
382	H O O	CN	×	
383	H 07 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	NO ₂	×<	\(\sigma^\circ\)

17. (Currently amended) The compound according to claim 16, wherein said compound is selected from compound
numbers: 54, 209, 237, 281, 295, 309, and 368, wherein the compound numbers are as defined below: [[.]]

	A	R ⁸	D'	E
209	H 0 1	s	×<	
237	H°, ,	~ King s	×	(T)
281	H° N	× H N N	×	() (1)

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295	H, O, J,	, In	×<	→ OH Me
309	H° N	-CONHMe	×	(C)
368	H O N	× S	×	\(\sqrt{\sq}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}

- 18. (Previously presented) A composition comprising a compound according to claim 1 or 9, in an amount sufficient to inhibit an aspartyl protease; and a pharmaceutically acceptable carrier.
- 19. (Original) The composition according to claim
 18, wherein said composition is in a pharmaceutically
 acceptable form for administration to a human being.
- 20. (Original) The composition according to claim 18, wherein said composition additionally comprises an additional anti-viral agent.
- 21. (Original) The composition according to claim
 18, wherein said composition comprises at least one additional
 therapeutic agent selected from (1 alpha, 2 beta, 3 alpha)-9[2,3-bis(hydroxymethyl)cyclobutyl]- guanine [(-)BHCG, SQ-

34514]; oxetanocin-G (3,4-bis-(hydroxymethyl)-2oxetanosyl]guanine); acyclic nucleosides, such as acyclovir, valaciclovir, famciclovir, ganciclovir or penciclovir; acyclic nucleoside phosphonates, such as (S)-1-(3-hydroxy-2phosphonyl-methoxypropyl)cytosine (HPMPC); ribonucleotide reductase inhibitors, such as 2-acetylpyridine 5-[(2chloroanilino)thiocarbonyl) thiocarbonohydrazone, 3'azido-3'deoxythymidine; other 2',3'-dideoxynucleosides such as 2',3'dideoxycytidine, 2',3'-dideoxyadenosine, 2',3'-dideoxyinosine, or 2',3'-didehydrothymidine; other aspartyl protease inhibitors, such as indinavir, ritonavir, nelfinavir, or [3S-[3R*(1R*, 2S*)]]-[3[[(4-aminophenyl)sulfonyl](2methylpropyl)amino]-2-hydroxy-1-(phenylmethyl)propyl]tetrahydro-3-furanyl ester (amprenavir); oxathiolane nucleoside analogues, such as (-)-cis-1-(2-hydroxymethyl)-1,3oxathiolane 5-yl)-cytosine (lamivudine) or cis-1-(2-(hydroxymethyl)-1,3-oxathiolan-5-yl)-5-fluorocytosine (FTC); 3'-deoxy-3'-fluorothymidine; 5-chloro-2',3'-dideoxy-3'fluorouridine; (-)-cis-4-[2-amino-6-(cyclopropylamino)-9Hpurin-9-yl]-2-cyclopentene-1-methanol; ribavirin; 9-[4hydroxy-2-(hydroxymethyl)but-1-yl]-guanine (H2G); tat inhibitors, such as 7-chloro-5-(2-pyrryl)-3H-1,4benzodiazepin-2-(H)one (Ro5-3335) or 7-chloro-1,3-dihydro-5-(1H-pyrrol-2yl)-3H-1,4-benzodiazepin-2-amine (Ro24-7429);

interferons, such as α-interferon; renal excretion inhibitors such as probenecid; nucleoside transport inhibitors such as dipyridamole; pentoxifylline; N-acetylcysteine (NAC); Procysteine; α -trichosanthin; phosphonoformic acid; immunomodulators, such as interleukin II or thymosin; granulocyte macrophage colony stimulating factors; erythropoetin; soluble CD4 and genetically engineered derivatives thereof; non-nucleoside reverse transcriptase inhibitors (NNRTIs), such as nevirapine (BI-RG-587), loviride (α -APA) or delavuridine (BHAP); phosphonoformic acid; 1,4-dihydro-2H-3,1-benzoxazin-2-ones NNRTIs, such as (-)-6-chloro-4-cyclopropylethynyl-4-trifluoromethyl-1,4-dihydro-2H-3,1-benzoxazin-2-one (L-743,726 or DMP-266); or quinoxaline NNRTIs, such as isopropyl (2S)-7-fluoro-3,4-dihydro-2-ethyl-3-oxo-1(2H)-quinoxalinecarboxylate (HBY1293).

- 22. (Previously presented) The composition according to claim 18, wherein said composition is in an orally available dosage form.
- 23. (Withdrawn) A method of treating a patient infected with a virus that depends upon an aspartyl protease for an obligatory event in its life cycle comprising the step

of administering to said patient a composition according to claim 18.

- 24. (Withdrawn) A method of treating a patient infected with HIV-I or HIV-II comprising the step of administering to said patient a composition according to claim 18.
- (Withdrawn) The method according to claim 23, 25. comprising the additional step of administering to said patient an additional therapeutic agent selected from (1 alpha, 2 beta, 3 alpha)-9-[2,3-bis(hydroxymethyl) cyclobutyl]quanine [(-)BHCG, SQ-34514]; oxetanocin-G (3,4-bis-(hydroxymethyl)-2-oxetanosyl]guanine); acyclic nucleosides, such as acyclovir, valaciclovir, famciclovir, ganciclovir or penciclovir; acyclic nucleoside phosphonates, such as (S)-1-(3-hydroxy-2-phosphonyl-methoxypropyl)cytosine (HPMPC); ribonucleotide reductase inhibitors, such as 2-acetylpyridine 5-[(2-chloroanilino)thiocarbonyl) thiocarbonohydrazone, 3'azido-3'-deoxythymidine; other 2',3'-dideoxynucleosides such as 2',3'-dideoxycytidine, 2',3'-dideoxyadenosine, 2',3'dideoxyinosine, or 2',3'-didehydrothymidine; other aspartyl protease inhibitors, such as indinavir, ritonavir, nelfinavir, or [3S-[3R*(1R*, 2S*)]]-[3[[(4-aminophenyl)sulfonyl](2-

methylpropyl)amino]-2-hydroxy-1-(phenylmethyl)propyl]tetrahydro-3-furanyl ester (amprenavir); oxathiolane nucleoside analogues, such as (-)-cis-1-(2-hydroxymethyl)-1,3oxathiolane 5-yl)-cytosine (lamivudine) or cis-1-(2-(hydroxymethyl)-1,3-oxathiolan-5-yl)-5-fluorocytosine (FTC); 3'-deoxy-3'-fluorothymidine; 5-chloro-2',3'-dideoxy-3'fluorouridine; (-)-cis-4-[2-amino-6-(cyclopropylamino)-9Hpurin-9-yl]-2-cyclopentene-1-methanol; ribavirin; 9-[4hydroxy-2-(hydroxymethyl)but-1-yl]-quanine (H2G); tat inhibitors, such as 7-chloro-5-(2-pyrryl)-3H-1,4benzodiazepin-2-(H)one (Ro5-3335) or 7-chloro-1,3-dihydro-5-(1H-pyrrol-2yl)-3H-1,4-benzodiazepin-2-amine (Ro24-7429); interferons, such as α -interferon; renal excretion inhibitors such as probenecid; nucleoside transport inhibitors such as dipyridamole; pentoxifylline; N-acetylcysteine (NAC); Procysteine; α -trichosanthin; phosphonoformic acid; immunomodulators, such as interleukin II or thymosin; granulocyte macrophage colony stimulating factors; erythropoetin; soluble CD4 and genetically engineered derivatives thereof; non-nucleoside reverse transcriptase inhibitors (NNRTIs), such as nevirapine (BI-RG-587), loviride (α -APA) or delavuridine (BHAP); phosphonoformic acid; 1,4dihydro-2H-3,1-benzoxazin-2-ones NNRTIs, such as (-)-6-chloro-

4-cyclopropylethynyl-4-trifluoromethyl-1,4-dihydro-2H-3,1-benzoxazin-2-one (L-743,726 or DMP-266); or quinoxaline NNRTIs, such as isopropyl (2S)-7-fluoro-3,4-dihydro-2-ethyl-3-oxo-1(2H)-quinoxalinecarboxylate (HBY1293), wherein said additional agent is administered to said patient as either a separate dosage form or as a single dosage form together with said compound.

- 26. (Withdrawn) A method of treating a patient diagnosed with AIDS; AIDS related complex (ARC); progressive generalized lymphadenopathy (PGL); Kaposi's sarcoma, thrombocytopenic purpura; AIDS-related neurological conditions such as AIDS dementia complex, multiple sclerosis or tropical paraperesis; anti-HIV antibody-positive conditions; or HIV-positive conditions, comprising the step of administering to said patient a composition according to claim 18.
- 27. (Withdrawn) The method according to claim 26, comprising the additional step of administering to said patient an additional therapeutic agent selected from (1 alpha, 2 beta, 3 alpha)-9-[2,3-bis(hydroxymethyl) cyclobutyl]guanine [(-)BHCG, SQ-34514]; oxetanocin-G (3,4-bis-(hydroxymethyl)-2-oxetanosyl]guanine); acyclic nucleosides, such as acyclovir, valaciclovir, famciclovir, ganciclovir or

penciclovir; acyclic nucleoside phosphonates, such as (S)-1-(3-hydroxy-2-phosphonyl-methoxypropyl)cytosine (HPMPC); ribonucleotide reductase inhibitors, such as 2-acetylpyridine 5-[(2-chloroanilino)thiocarbonyl) thiocarbonohydrazone, 3'azido-3'-deoxythymidine; other 2',3'-dideoxynucleosides such as 2',3'-dideoxycytidine, 2',3'-dideoxyadenosine, 2',3'dideoxyinosine, or 2',3'-didehydrothymidine; other aspartyl protease inhibitors, such as indinavir, ritonavir, nelfinavir, or [3S-[3R*(1R*, 2S*)]]-[3[[(4-aminophenyl)sulfonyl](2methylpropy1)amino]-2-hydroxy-1-(phenylmethyl)propy1]tetrahydro-3-furanyl ester (amprenavir); oxathiolane nucleoside analogues, such as (-)-cis-1-(2-hydroxymethyl)-1,3oxathiolane 5-yl)-cytosine (lamivudine) or cis-1-(2-(hydroxymethyl)-1,3-oxathiolan-5-yl)-5-fluorocytosine (FTC); 3'-deoxy-3'-fluorothymidine; 5-chloro-2',3'-dideoxy-3'fluorouridine; (-)-cis-4-[2-amino-6-(cyclopropylamino)-9Hpurin-9-yl]-2-cyclopentene-1-methanol; ribavirin; 9-[4hydroxy-2-(hydroxymethyl)but-1-yl]-guanine (H2G); tat inhibitors, such as 7-chloro-5-(2-pyrryl)-3H-1,4benzodiazepin-2-(H)one (Ro5-3335) or 7-chloro-1,3-dihydro-5-(1H-pyrrol-2yl)-3H-1,4-benzodiazepin-2-amine (Ro24-7429); interferons, such as α -interferon; renal excretion inhibitors such as probenecid; nucleoside transport inhibitors such as

dipyridamole; pentoxifylline; N-acetylcysteine (NAC); Procysteine; α -trichosanthin; phosphonoformic acid; immunomodulators, such as interleukin II or thymosin; granulocyte macrophage colony stimulating factors; erythropoetin; soluble CD4 and genetically engineered derivatives thereof; non-nucleoside reverse transcriptase inhibitors (NNRTIs), such as nevirapine (BI-RG-587), loviride $(\alpha - APA)$ or delavuridine (BHAP); phosphonoformic acid; 1,4dihydro-2H-3,1-benzoxazin-2-ones NNRTIs, such as (-)-6-chloro-4-cyclopropylethynyl-4-trifluoromethyl-1,4-dihydro-2H-3,1benzoxazin-2-one (L-743,726 or DMP-266); or quinoxaline NNRTIs, such as isopropyl (2S)-7-fluoro-3,4-dihydro-2-ethyl-3oxo-1(2H)-quinoxalinecarboxylate (HBY1293), wherein said additional agent is administered to said patient as either a separate dosage form or as a single dosage form together with said compound.

28. (Currently amended) The compound according to claim 15, wherein said compound is compound number 368.

wherein said compound number is as defined below: [[.]]

	A	R ⁸	D'	Е
368	H°,	, N	×	

- 29. (Previously presented) The composition according to claim 19, wherein said composition is in an orally available dosage form.
- 30. (Previously presented) The composition according to claim 20, wherein said composition is in an orally available dosage form.
- 31. (Previously presented) The composition according to claim 21, wherein said composition is in an orally available dosage form.
- 32. (Withdrawn) The method according to claim 24, comprising the additional step of administering to said

patient an additional therapeutic agent selected from (1 alpha, 2 beta, 3 alpha)-9-[2,3-bis(hydroxymethyl) cyclobutyl]quanine [(-)BHCG, SQ-34514]; oxetanocin-G (3,4-bis-(hydroxymethyl)-2-oxetanosyl]guanine); acyclic nucleosides, such as acyclovir, valaciclovir, famciclovir, ganciclovir or penciclovir; acyclic nucleoside phosphonates, such as (S)-1-(3-hydroxy-2-phosphonyl-methoxypropyl)cytosine (HPMPC); ribonucleotide reductase inhibitors, such as 2-acetylpyridine 5-[(2-chloroanilino)thiocarbonyl) thiocarbonohydrazone, 3'azido-3'-deoxythymidine; other 2',3'-dideoxynucleosides such as 2',3'-dideoxycytidine, 2',3'-dideoxyadenosine, 2',3'dideoxyinosine, or 2',3'-didehydrothymidine; other aspartyl protease inhibitors, such as indinavir, ritonavir, nelfinavir, or [3S-[3R*(1R*, 2S*)]]-[3[[(4-aminophenyl)sulfonyl](2methylpropyl)amino]-2-hydroxy-1-(phenylmethyl)propyl]tetrahydro-3-furanyl ester (amprenavir); oxathiolane nucleoside analogues, such as (-)-cis-1-(2-hydroxymethyl)-1,3oxathiolane 5-yl)-cytosine (lamivudine) or cis-1-(2-(hydroxymethyl)-1,3-oxathiolan-5-yl)-5-fluorocytosine (FTC); 3'-deoxy-3'-fluorothymidine; 5-chloro-2',3'-dideoxy-3'fluorouridine; (-)-cis-4-[2-amino-6-(cyclopropylamino)-9Hpurin-9-yl]-2-cyclopentene-1-methanol; ribavirin; 9-[4hydroxy-2-(hydroxymethyl)but-1-yl]-quanine (H2G); tat inhibitors, such as 7-chloro-5-(2-pyrryl)-3H-1,4-

benzodiazepin-2-(H)one (Ro5-3335) or 7-chloro-1,3-dihydro-5-(1H-pyrrol-2yl)-3H-1,4-benzodiazepin-2-amine (Ro24-7429); interferons, such as α -interferon; renal excretion inhibitors such as probenecid; nucleoside transport inhibitors such as dipyridamole; pentoxifylline; N-acetylcysteine (NAC); Procysteine; α -trichosanthin; phosphonoformic acid; immunomodulators, such as interleukin II or thymosin; granulocyte macrophage colony stimulating factors; erythropoetin; soluble CD4 and genetically engineered derivatives thereof; non-nucleoside reverse transcriptase inhibitors (NNRTIs), such as nevirapine (BI-RG-587), loviride $(\alpha - APA)$ or delavuridine (BHAP); phosphonoformic acid; 1,4dihydro-2H-3,1-benzoxazin-2-ones NNRTIs, such as (-)-6-chloro-4-cyclopropylethynyl-4-trifluoromethyl-1,4-dihydro-2H-3,1benzoxazin-2-one (L-743,726 or DMP-266); or quinoxaline NNRTIs, such as isopropyl (2S)-7-fluoro-3,4-dihydro-2-ethyl-3oxo-1(2H)-quinoxalinecarboxylate (HBY1293), wherein said additional agent is administered to said patient as either a separate dosage form or as a single dosage form together with said compound.

33. (Currently amended) A compound selected from compound numbers: 26, 27, 28, 29, 30, 31, 32, 32, 33, 34, 35,

36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 50, 51, 52, 53, 54, 55, 56, 57, 59, 60, 61, 62, 63, 71, 72, 73, 74, 75, 201, 202, 203, 204, 205, 206, 207, 208, 209, 210, 211, 212, 213, 214, 215, 216, 217, 218, 219, 220, 221, 222, 223, 224, 225, 226, 227, 228, 229, 230, 231, 232, 233, 234, 235, 236, 237, 238, 239, 240, 241, 242, 243, 244, 245, 246, 247, 248, 249, 250, 251, 252, 253, 254, 255, 256, 257, 258, 259, 260, 261, 262, 263, 264, 265, 266, 267, 268, 269, 270, 271, 272, 273, 274, 275, 276, 277, 278, 279, 280, 281, 282, 283, 284, 285, 286, 287, 288, 289, 290, 291, 292, 293, 294, 295, 296, 297, 298, 299, 300, 301, 302, 303, 304, 305, 309, 310, 311, 312, 313, 314, 315, 316, 317, 318, 319, 320, 321, 322, 323, 324, 326, 327, 329, 331, 332, 333, 334, 335, 336, 337, 338, 339, 340, 341, 342, 343, 344, 345, 346, 347, 348, 349, 351, 352, 353, 354, 355, 356, 357, 358, 359, 361, 364, 365, 366, 367, 368, 369, 370, 371, 372, 373, 374, 375, 376, 377, 378, 379, 380, 381, 382, 383, 384, 385, 386, 387, 388, 389, 390, 391, 392, 393, 394, 395, 396, 397 and 398, wherein said compound numbers are as defined below: [[.]]

	A	S(O) ₂ E	R ⁸
26			
26	H H	NH ₂	
27	H H	NH ₂	
28	H H	NH ₂	
29	H H	NH ₂	
30	H H	NH ₂	
31	H H	NH ₂	N

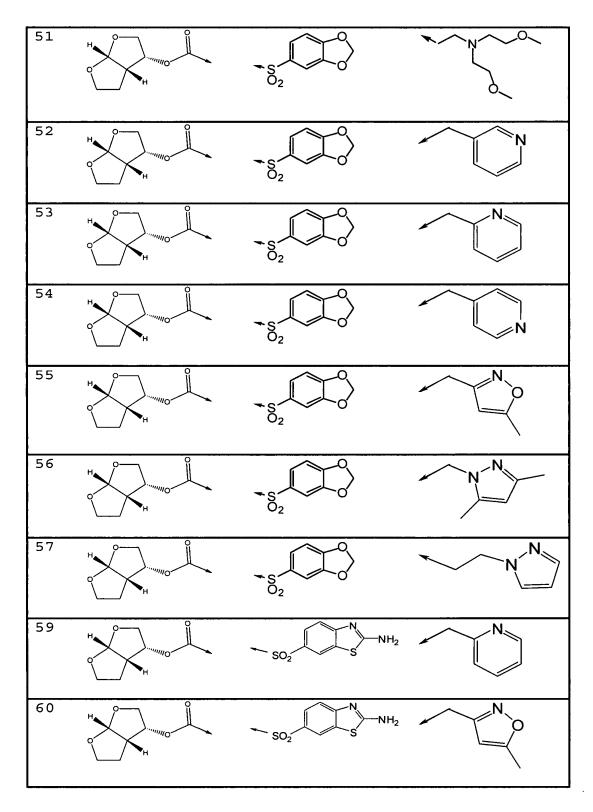
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32	H	NH ₂	Н
33	H H	S _{O₂} NH ₂	N
34	H H	NH ₂	
35	H	S NH ₂	N
36	H H	S NH ₂	N
37	H H	S NH ₂	
38	H	SO ₂ OH	
39	H	SO ₂ OH	Н
40	H	SO ₂ OH	

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41	H H	SO ₂ OH	
42	H H	SO ₂ OH	
43	H H	SO ₂ OH	N
44	H O H	SO ₂ NH ₂	
46	H H	- S ₂	
47	H O H	- §	Н
48	H 0 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	- §	
49	H H	- S ₂	NO
50	H H	-S ₂	, L

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A
$$S(0)_2E$$
 R^8

71 NH_2 NH_2

72 NH_2 NH

	A	S(O) ₂ E	R ⁸
73	H H	S ₂	N
74	H H	S NH ₂	
75	H	S NH ₂	

	A	R ⁸	D'	E
201	H, O, J, O, J, O,	-Н	×	(T)
202	H O O		×<	\(\sigma^{\circ}\)
203	H O N	, O	×	XX°

204	но о	CF ₃	/	△ -0
	ON HOUSE	<i>;</i> ~~	×. <	
205	H O N	× JF	\times	
206	H ° ,	OMe	×	XX°
207	H° N	↓\bigcip\s\	х-<	\(\sigma\)
208	H O J	× S	×	XX°
209	H, O, J,	× S	×<	\(\sigma_{\circ}\)
210	H° N	, L3	×<	(C)
211	H°,	×	х. <	\(\frac{1}{2}\)
212	H° N	XIN	×	(T)
213	H, O, J,	S	×	↓Ç;
214	H, O, J,	i Ch	×. <	
215	H 0 1	× CN	×. <	\(\sigma^{\circ}\)
216	H O J	CF,	×<	\(\sigma_{\circ}^{\circ}\)

			 	
217	H O N	` ,~~ CF ₃	×<	\(\cdot\)
218	H O N	NH ₂	\times	
219	H ° N	NHCOOt-bu	×	
220	H, O, J,	NH ₂	×<	\(\sqrt{\sq}\sqrt{\sq}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}}
221	H°,	NHCOOt-bu	×<	(T)
222	H°, Å	`NH₂	х. <	\(\sigma\)
223	H ° N	NHCOO-t-bu	×<	,(C);
224	H°,	, NH ₂	×<	\(\sigma\)
225	H ° ,	NHAC	\times	XX;
226	H ° N	NHCOOMe	×<	XX;
227	H°,	NHSO ₂ Me	\times	
228	H°,	NHCONHMe	х. <	
229	H O N	NHAc	×<	XX°
230	H O O	NHAC	×<	

231	H° N	X P	×.<	\(\sigma\)
232	H° N	××T	×<	(X)
233	H, O, J, O,	× N°	×<	(X)
234	H° N	÷ N	×<	\(\sigma\)
235	H° N	× Nº CO	×<	(X)
236	H O N		×	
237	H° N	~ H S	×<	\(\sigma\)
238	H ° N	× Y S	×<	
239	HO HO	i s	×<	
240	H, O, J,	OMe	×<	
241	H, O, J, O,	N OMe	×	
242	H, O, I, O,	, H O OMe	×	XX°
243	H° N	N OMe	×	XX°
244	H, O, J,	H O OMe	×<	\(\mathcal{L}\)
	·			

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0.45				
245		N OEt	×. <	
246	H, O, J,	N OE	×. <	
247	H ° ,	, N YO OEt	×	
248	H ° ,	N O-i-Pr	×<	
249	H ° ,	N O-i-Pr	×<	
250	H ° ,	HO O-i-Pr	×<	XX;
251	H ° ,	NHSO ₂ Me	×	XX;
252	H°,	, NHSO₂Me	×	\(\sigma\)
253	H°,	NHCONMe ₂	×	\(\sigma\)
254	H ° ,	`NHCONMe₂	×. <	XX;
255	H, O, J,	NHCONMe ₂	×. <	
256	H 0 1 0 1 0 1 0 1 0 1 0 1 0 1 0 1 0 1 0	NHCONHMe	×<	
257	H, O, J,	, NHCONHMe	×<	
258	H O J		×<	

259	H O N	NHCSNHMe	×<	(C)
260	H ° ,	, NHCSNHMe	\times	
261	H ° N	× H LO	×<	\(\sigma\)
262	H ° ,	N.CN NOPh	х. <	XX°
263	H ° N	N.CN N.OPh	×.<	XX;
264	H O O	N.CN N NH ₂	×.<	XX°
265	H O O	N-CN NHMe	×<	XX°
266	H O H	+ H N O	×<	XX°
267	H ° ,	N.CN N NH ₂	×<	II.
268	H ° ,	N.CN N NHMe	×. <	XX°
269	H ° N	N.CN N.Me ₂	×<	\(\sigma\)
270	H O N	× N N CN	×<	\(\sigma_{\circ}\)
271	HO	N N N	×.<	XX°>

272	H O	N CEL ³	×	
273	H O O	NMe ₂	\times	
274	H ° À	NMe ₂	×<	\(\sigma^\circ\)
275	H ° ,	NMe ₂	х-<	(C)
276	H ° ,	*~1	х-<	(C)
277	H ° N	X_N_S	х-<	XX°
278	H°,	× NON	×	(T)
279	H O J	- N.N.	×<	\(\sigma_{\circ}\)
280	H, O, J,	×~~	×<	(T)
281	H° N	× H NN	×<	\(\sigma\)
282	H°,	NNMe	×. <	\(\)
283	H° N	OMe	×	(T)
284	H ° , , , , , , , , , , , , , , , , , ,	X_N	×	(C)
285	H ° 7, ° 1, ° 1, ° 1, ° 1, ° 1, ° 1, ° 1,	CONH ₂	×	\(\sigma\)

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286	H, O, J,	CONHMe	×	\(\sigma\)
287	H, O, J, O,	CONMe ₂	×	\(\sigma\)
288	H ° ,	CONHBU	×<	XX.
289	H O N	; O-cn	×	NH ₂
290	H O J	;D	×<	NH ₂
291	H, O, J, O,	Н	×<	NH ₂
292	H°,	NH ₂	×	NH ₂
293	H°,	; In	×	NH ₂
294	H O J	, SN	×<	OAc Me
295	H O J	; SN	×<	OH Me
296	H, O, J,	<i>,</i> ,0	×<	NH ₂
297	H O O	O, SHO	×<	NO ₂
298	H, O, J,	ON NH2	×. <	NH ₂

299	HOTHO	; In	×<	NH ₂
300	H° N	, S N	×. <	NH ₂
301	H°,	, D-cn	х-<	NH ₂
302	H	, D-cn	×<	, oh
303	H° N	COOMe	×	
304	H O N	√ CF ₃	×	XX°>
305	H O N		,×	
309	H° N	-CONHMe	×.<	XX;
310	H° N	-CONH-i-Pr	×	XX;
311	H° N	-CONMe2	х. <	XX°
312	H ° ,	-CONH2	×<	XX°
313	H ° ,	Y N N N	×<	XX°
314	H, O, J,	CONH ₂	х-<	XX°
315	H O N	CONHMe	×	

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316	H ° ,	`OSO₃K	×<	
317	H ° N	,,,, он	×<	XX°
318	H ° À	OCONHMe	×<	\(\tau^\chi_\)
319	H ° N	УОН	×<	
320	HOTH	OCONHMe	×<	
321	H ° ,	OCONH ₂	×.<	\(\sigma\)
322	H 0 1	; ⁄-он	×-<	\(\sigma\)
323	H ° ,	OCONHMe	×<	\(\frac{1}{2}\)
324	H°,	*/ N S	×<	\(\tau^\chi_\)
326	H O H	СООМе	×<	\(\mathcal{C}\)
327	H ° N	>∕_соон	×<	\(\sigma\)
329	H O H	,O	√ ₀₊	
331	H O J	NH NMe₂	×<	
332	H O O	FF	×<	

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333	H°,	021	\times	
334	H, O, A,	O ₂ N COOH	×.<	
335	H°, ,	NC (N	×	XX°
336	H, O, J,	O ₂ N NO ₂	×<	
337	H°,	÷ [n	х. <	
338	H O J	Dcn	×<	
339	H°,	СООН	х. <	\(\frac{1}{2}\)
340	H°,		\times	√ V _F
341	H°,	Н	×<	√ V _F
342	HOTHO	CN	×.<	XF OXF
343	H°,	,0	×.<	OAc
344	H°,	÷O.	×<	OH
345	H°,	Н	×<	OAc

346	H O O	CN	\times	OH
347	H°,	, , , , s	\times	OAc
348	H, O, J, O,	× S _N	×. <	↓ OH
349	H, O, J,	, , , , , , , , , , , , , , , , , , ,	×.<	
351	H°,	Н	×. <	(T)
352	H° N	× O	×. <	(T)
353	H, O, J,	<i>†</i>	×. <	XX°
354	H ° N	NH ₂	×.<	\(\sigma\)
355	H, O, J, O,	NHAC	х. <	
356	H° N	NHCOOMe	×<	\(\sigma\)
357	H° N	, NHSO₂Me	×<	XX°
358	H°, i	, NHCONHMe	×<	\(\sigma\)
359	H, O, J,	NMe ₂	×<	
361	H°,		OCONHM	XX°

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			•	
364	H O J		×<	
365	H ° ,	,D	×<	
366	H, O, J,	÷	×<	
367	H O J	; Dcn	×<	
368	H° N	, S	×. <	XX.
369	H O J	SMe	×<	\(\sigma_{\circ}^{\circ}\)
370	H O N	, N-0	×	\(\sigma_{\circ}^{\circ}\)
371	H O J	F	×. <	XX°>
372	H O J	CF ₃	×<	
373	H O N	CF ₃	×	
374	H, O, J,	Sin	×<	\(\infty\)?
375	H O J	N-0	×. <	

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376	H O H		×<	\(\frac{1}{2}\)
377	H O N		×<	(T)
378	H ° ,	, DF	×<	(T)
379	H O I	F	×. <	XX°>
380	H O J	F	×. <	\(\sigma_{\circ}\)
381	HON	CN	×<	XX°>
382	H O N	CN	×	XX°
383	HOTA	NO ₂	×. <	XX°>
384	H O J	NO ₂	×	\(\sigma_{\circ}\)
385	H ° ,	+ Con	×. <	XX°
386	H°,	CI S, N	×<	(C)
387	H, O, J,	Š	×. <	\(\sigma_{\circ}\)
388	HOTHO	;↓ SO₂Ph	×	\(\sigma_{\circ}\)

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389	H O J	Z Z Z Z Z Z Z Z Z Z Z Z Z Z Z Z Z Z Z	×<	
390	H O O	OMe	×<	
391	H O J	OMe	×<	
392	H ° N	NC T	×	\(\sigma^\circ\)
393	H O N	· }.	\times	(XX)
394	H O O		×-<	
395	H O N	; O	×	XX°
396	H O H	CN	×O	\(\sigma_{\circ}^{\circ}\)
397	H°,	∤ CN	×	II.
398	H, O, J,	, In	×	\(\sigma_{\circ}\)